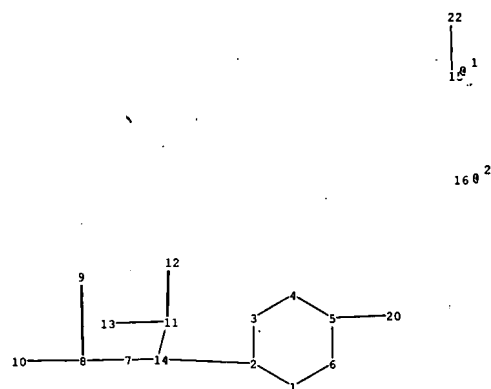
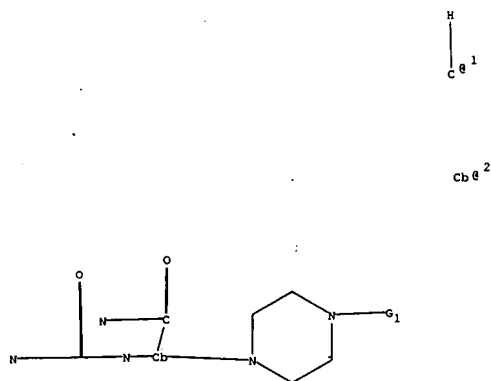


EAST Search History

Ref #	Hits	Search Query	DBs	Default Operator	Plurals	Time Stamp
L4	3318	((544/357) or (544/377) or (544/381) or (544/396) or (544/400) or (540/575)).CCLS.	US-PGPUB; USPAT	OR	OFF	2007/04/28 18:35
L5	48	1-piperazinyl near25 benzamide	US-PGPUB; USPAT	OR	OFF	2007/04/28 18:36
L6	13	I4 and I5	US-PGPUB; USPAT	OR	OFF	2007/04/28 18:37
L7	35	I5 not I6	US-PGPUB; USPAT	OR	OFF	2007/04/28 18:37



chain nodes :

7 8 9 10 11 12 14 15 16 20 22

ring nodes :

1 2 3 4 5 6

ring/chain nodes :

13

chain bonds :

2-14 5-20 7-8 7-14 8-9 8-10 11-12 11-13 11-14 15-22

ring bonds :

1-2 1-6 2-3 3-4 4-5 5-6

exact/norm bonds :

1-2 1-6 2-3 3-4 4-5 5-6 5-20 7-8 8-9 8-10 11-12 11-13

exact bonds :

2-14 7-14 11-14 15-22

isolated ring systems :

containing 1 :

G1:[*1],[*2]

Match level :

1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:CLASS8:CLASS9:CLASS10:CLASS11:CLASS

12:CLASS13:CLASS14:Atom 15:CLASS16:Atom 20:CLASS22:CLASS

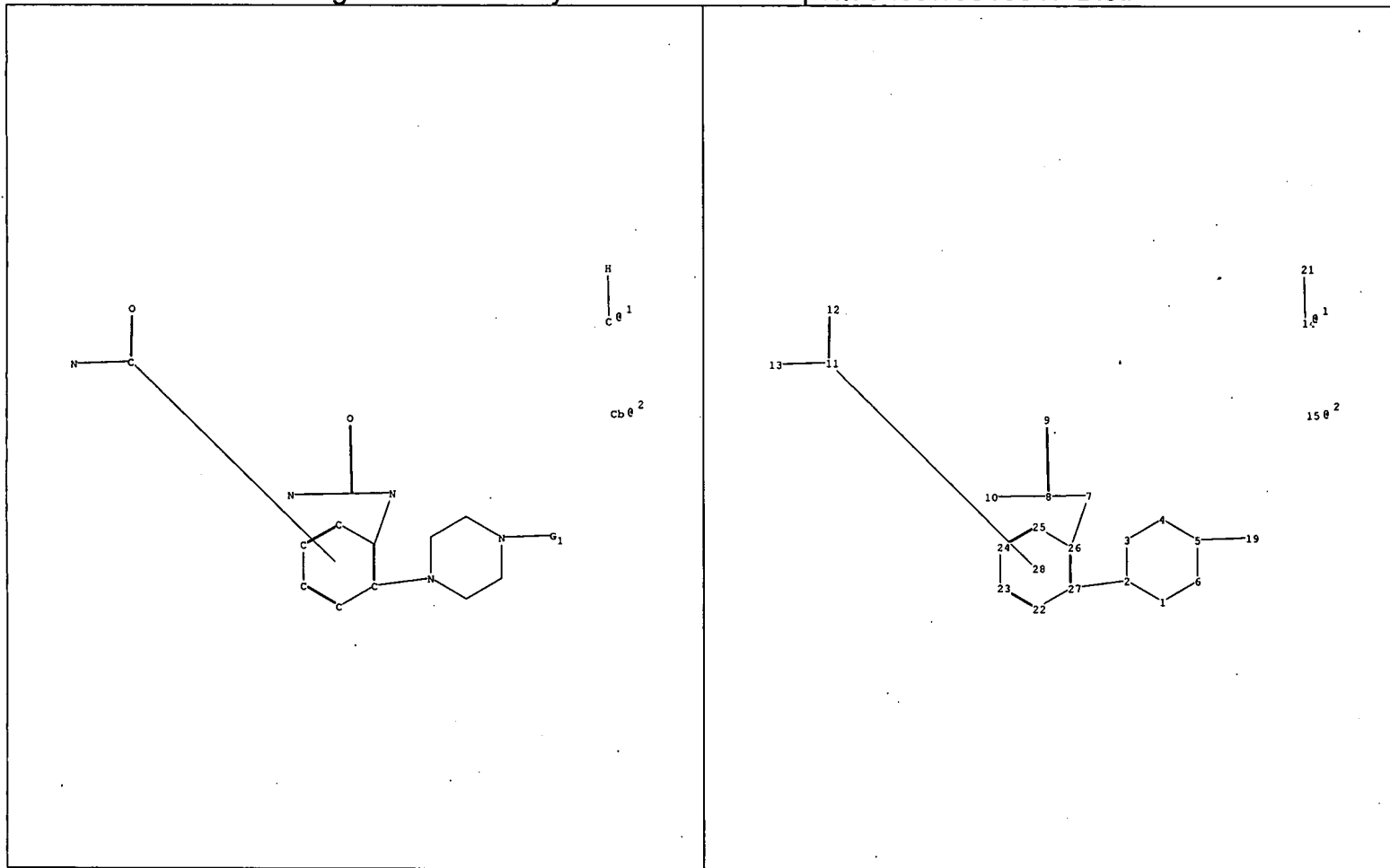
Generic attributes :

14:

Saturation : Unsaturated

Number of Carbon Atoms : less than 7

Type of Ring System : Monocyclic



chain nodes :

7 8 9 10 11 12 14 15 19 21

ring nodes :

1 2 3 4 5 6 22 23 24 25 26 27

ring/chain nodes :

13

chain bonds :

2-27 5-19 7-8 7-26 8-9 8-10 11-12 11-13 14-21

ring bonds :

1-2 1-6 2-3 3-4 4-5 5-6 22-23 22-27 23-24 24-25 25-26 26-27

exact/norm bonds :

1-2 1-6 2-3 2-27 3-4 4-5 5-6 5-19 7-8 7-26 8-9 8-10 11-12 11-13

exact bonds :

14-21

normalized bonds :

22-23 22-27 23-24 24-25 25-26 26-27

isolated ring systems :

containing 1 :

G1:[*1],[*2]

Match level :

1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:CLASS8:CLASS9:CLASS10:CLASS11:CLASS
12:CLASS

13:CLASS14:CLASS15:Atom 19:CLASS21:CLASS22:CLASS23:Atom 24:Atom 25:Atom
26:CLASS27:Atom 28:Atom

10/815017

=> s 11

SAMPLE SEARCH INITIATED 16:08:07 FILE 'REGISTRY'
SAMPLE SCREEN SEARCH COMPLETED - 1916 TO ITERATE

100.0% PROCESSED 1916 ITERATIONS
SEARCH TIME: 00.00.01

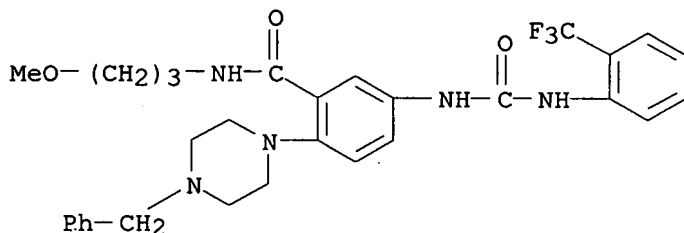
34 ANSWERS

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
BATCH **COMPLETE**
PROJECTED ITERATIONS: 35695 TO 40945
PROJECTED ANSWERS: 331 TO 1029

L2 34 SEA SSS SAM L1

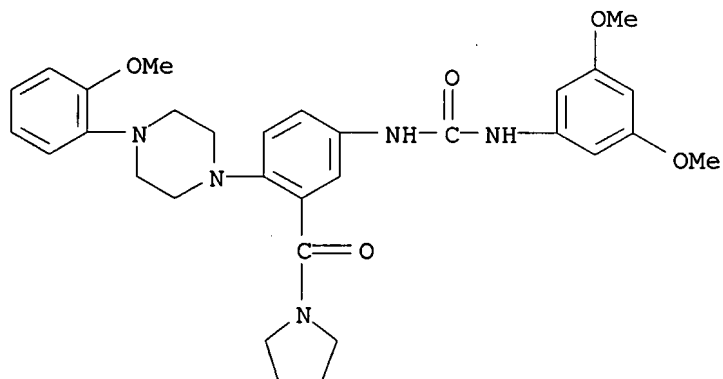
=> d 12 1 5 10 30

L2 ANSWER 1 OF 34 REGISTRY COPYRIGHT 2007 ACS on STN
RN 861982-63-0 REGISTRY
ED Entered STN: 29 Aug 2005
CN Benzamide, N-(3-methoxypropyl)-2-[4-(phenylmethyl)-1-piperazinyl]-5-[[[2-(trifluoromethyl)phenyl]amino]carbonyl]amino]- (9CI) (CA INDEX NAME)
MF C30 H34 F3 N5 O3
SR Chemical Library
Supplier: ComGenex International Inc.
LC STN Files: CHEMCATS



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

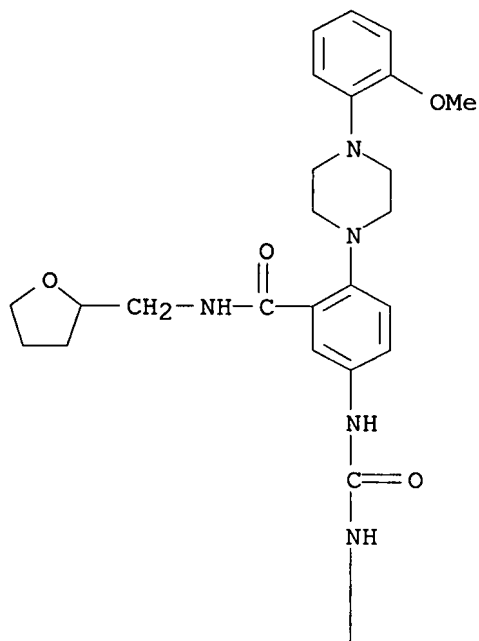
L2 ANSWER 5 OF 34 REGISTRY COPYRIGHT 2007 ACS on STN
RN 860078-73-5 REGISTRY
ED Entered STN: 12 Aug 2005
CN Pyrrolidine, 1-[5-[[[(3,5-dimethoxyphenyl)amino]carbonyl]amino]-2-[4-(2-methoxyphenyl)-1-piperazinyl]benzoyl]- (9CI) (CA INDEX NAME)
MF C31 H37 N5 O5
SR Chemical Library



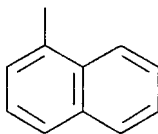
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 ANSWER 10 OF 34 REGISTRY COPYRIGHT 2007 ACS on STN
 RN 857668-00-9 REGISTRY
 ED Entered STN: 29 Jul 2005
 CN Benzamide, 2-[4-(2-methoxyphenyl)-1-piperazinyl]-5-[[1-(naphthalenylamino)carbonyl]amino]-N-[(tetrahydro-2-furanyl)methyl]- (9CI)
 (CA INDEX NAME)
 MF C34 H37 N5 O4
 SR Chemical Library
 Supplier: ComGenex International Inc.
 LC STN Files: CHEMCATS

PAGE 1-A



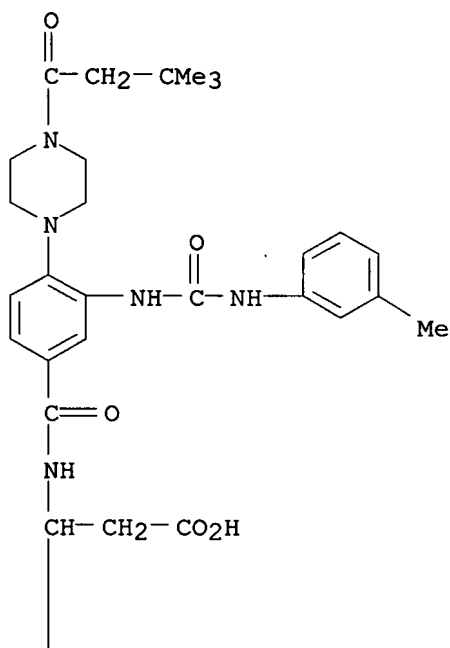
PAGE 2-A

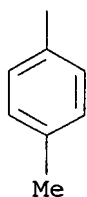


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 ANSWER 30 OF 34 REGISTRY COPYRIGHT 2007 ACS on STN
 RN 439274-60-9 REGISTRY
 ED Entered STN: 18 Jul 2002
 CN Benzenepropanoic acid, β -[[4-[4-(3,3-dimethyl-1-oxobutyl)-1-piperazinyl]-3-[[[(3-methylphenyl)amino]carbonyl]amino]benzoyl]amino]-4-methyl- (9CI) (CA INDEX NAME)
 MF C35 H43 N5 O5
 SR Chemical Library
 Supplier: Ambinter

PAGE 1-A





PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

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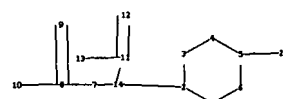
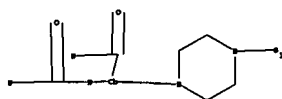
Uploading C:\Documents and Settings\EBernhardt\My Documents\Stnexp\Queries\10815017-A.str



1



2



chain nodes :
 7 8 9 10 11 12 14 15 16 20 22
 ring nodes :
 1 2 3 4 5 6
 ring/chain nodes :
 13

10/815017

chain bonds :
2-14 5-20 7-8 7-14 8-9 8-10 11-12 11-13 11-14 15-22
ring bonds :
1-2 1-6 2-3 3-4 4-5 5-6
exact/norm bonds :
1-2 1-6 2-3 3-4 4-5 5-6 5-20 7-8 8-9 8-10 11-12 11-13
exact bonds :
2-14 7-14 11-14 15-22
isolated ring systems :
containing 1 :

G1:[*1],[*2]

Match level :
1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:CLASS 8:CLASS 9:CLASS 10:CLASS
11:CLASS 12:CLASS 13:CLASS 14:Atom 15:CLASS 16:Atom 20:CLASS 22:CLASS
Generic attributes :
14:
Saturation : Unsaturated
Number of Carbon Atoms : less than 7
Type of Ring System : Monocyclic

L3 STRUCTURE UPLOADED

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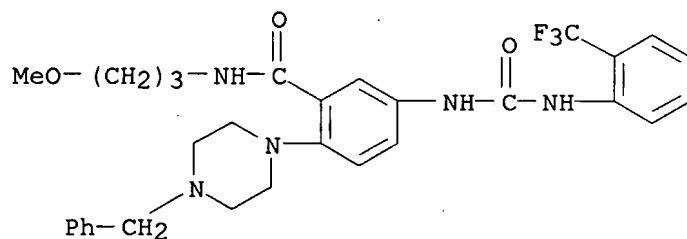
100.0% PROCESSED 1916 ITERATIONS 27 ANSWERS
SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
BATCH **COMPLETE**
PROJECTED ITERATIONS: 35695 TO 40945
PROJECTED ANSWERS: 229 TO 851

L4 27 SEA SSS SAM L3

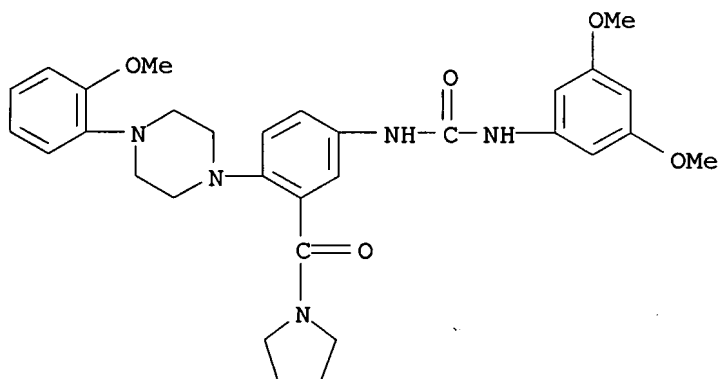
=> d 14 1 5 10 27

L4 ANSWER 1 OF 27 REGISTRY COPYRIGHT 2007 ACS on STN
RN 861982-63-0 REGISTRY
ED Entered STN: 29 Aug 2005
CN Benzamide, N-(3-methoxypropyl)-2-[4-(phenylmethyl)-1-piperazinyl]-5-[[[2-(trifluoromethyl)phenyl]amino]carbonyl]amino]- (9CI) (CA INDEX NAME)
MF C30 H34 F3 N5 O3
SR Chemical Library
Supplier: ComGenex International Inc.
LC STN Files: CHEMCATS



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

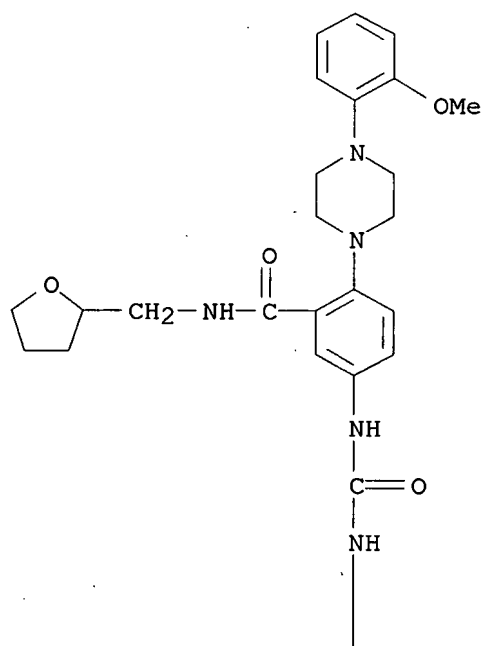
L4 ANSWER 5 OF 27 REGISTRY COPYRIGHT 2007 ACS on STN
 RN 860078-73-5 REGISTRY
 ED Entered STN: 12 Aug 2005
 CN Pyrrolidine, 1-[5-[[[(3,5-dimethoxyphenyl)amino]carbonyl]amino]-2-[4-(2-methoxyphenyl)-1-piperazinyl]benzoyl]- (9CI) (CA INDEX NAME)
 MF C31 H37 N5 O5
 SR Chemical Library



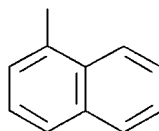
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L4 ANSWER 10 OF 27 REGISTRY COPYRIGHT 2007 ACS on STN
 RN 857668-00-9 REGISTRY
 ED Entered STN: 29 Jul 2005
 CN Benzamide, 2-[4-(2-methoxyphenyl)-1-piperazinyl]-5-[[[(1-naphthalenylamino)carbonyl]amino]-N-[(tetrahydro-2-furanyl)methyl]- (9CI) (CA INDEX NAME)
 MF C34 H37 N5 O4
 SR Chemical Library
 Supplier: ComGenex International Inc.
 LC STN Files: CHEMCATS

PAGE 1-A

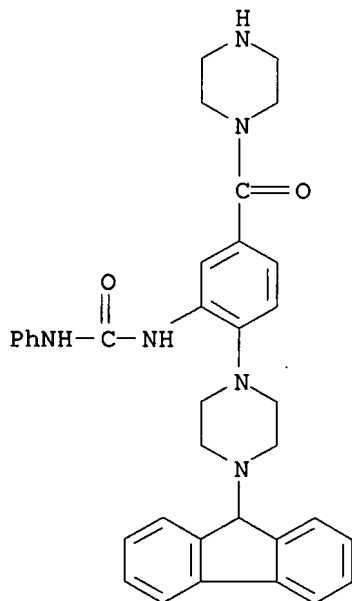


PAGE 2-A



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L4 ANSWER 27 OF 27 REGISTRY COPYRIGHT 2007 ACS on STN
 RN 773882-48-7 REGISTRY
 ED Entered STN: 02 Nov 2004
 CN Piperazine, 1-[4-[4-(9H-fluoren-9-yl)-1-piperazinyl]-3-
 [[(phenylamino)carbonyl]amino]benzoyl]- (9CI) (CA INDEX NAME)
 MF C35 H36 N6 O2
 SR CA
 LC STN Files: CA, CAPLUS, USPATFULL



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

1 REFERENCES IN FILE CA (1907 TO DATE)
1 REFERENCES IN FILE CAPLUS (1907 TO DATE)

=> s 13 sss full
FULL SEARCH INITIATED 16:12:06 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED - 39190 TO ITERATE

100.0% PROCESSED 39190 ITERATIONS 422 ANSWERS
SEARCH TIME: 00.00.02

L5 422 SEA SSS FUL L3

=> save l5
ENTER NAME OR (END):ten815017/A
ANSWER SET L5 HAS BEEN SAVED AS 'TEN815017/A'

=> file caplus		
COST IN U.S. DOLLARS	SINCE FILE	TOTAL
	ENTRY	SESSION
FULL ESTIMATED COST	190.85	191.06

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FILE COVERS 1907 - 28 Apr 2007 VOL 146 ISS 19

FILE LAST UPDATED: 27 Apr 2007 (20070427/ED)

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=> s 15

L6 1 L5

=> d 16 1 bib abs fhitr

L6 ANSWER 1 OF 1 CAPLUS COPYRIGHT 2007 ACS on STN

AN 2004:857583 CAPLUS

DN 141:332220

TI A preparation of (piperazinyphenyl)urea derivatives as phospholipase C
inhibitors, useful for the treatment of inflammatory disorders

IN Lagu, Bharat; Wachter, Michael; Rupert, Kenneth; Wachter, Michael

PA Janssen Pharmaceutica N.V., Belg.

SO PCT Int. Appl., 141 pp.

CODEN: PIXXD2

DT Patent

LA English

FAN.CNT 1

	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
PI	WO 2004087685	A2	20041014	WO 2004-US9846	20040331
	WO 2004087685	A3	20041216		

W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BW, BY, BZ, CA, CH,
CN, CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, EG, ES, FI, GB, GD,
GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC,
LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NA, NI,
NO, NZ, OM, PG, PH, PL, PT, RO, RU, SC, SD, SE, SG, SK, SL, SY,
TJ, TM, TN, TR, TT, TZ, UA, UG, US, UZ, VC, VN, YU, ZA, ZM, ZW,
RW: BW, GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZM, ZW, AM, AZ,
BY, KG, KZ, MD, RU, TJ, TM, AT, BE, BG, CH, CY, CZ, DE, DK, EE,
ES, FI, FR, GB, GR, HU, IE, IT, LU, MC, NL, PL, PT, RO, SE, SI,
SK, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN,
TD, TG

	US 2004235855	A1	20041125	US 2004-815017	20040331
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PRAI	US 2003-458938P	P	20030331		
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OS MARPAT 141:332220

GI

* STRUCTURE DIAGRAM TOO LARGE FOR DISPLAY - AVAILABLE VIA OFFLINE PRINT *

AB The invention relates to a preparation of (piperazinyphenyl)urea derivs. of

formula I [wherein: X is NH₂, NH-alkyl, NHOH, NH-CN, or heterocyclic ring, etc.; Y is one or more (un)substituted alkyl; Z is (CH₂)₂₋₅; R₁ is (un)substituted alkyl, cycloalkyl, or aryl, etc.; R₂ is (un)substituted alkyl, C(O)alkyl, C(O)alkenyl, aryl, or cycloalkyl, etc.; R₃ is O or S], useful as PLC-β₂ inhibitors. For instance, (piperazinyphenyl)urea derivative II (IC₅₀ = 1.2 μM) was prepared via addition of resin-bound (piperazinyphenyl)amine derivative III to Ph-N=C=O and subsequent resin cleavage (example 1).

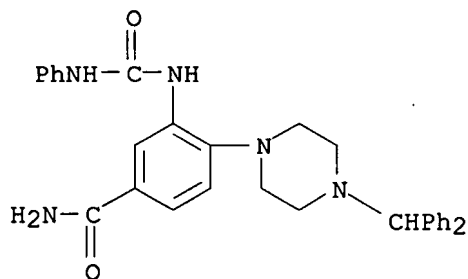
IT 773882-10-3P

RL: PAC (Pharmacological activity); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)

(preparation of (piperazinyphenyl)urea derivs. useful as PLC-β₂ inhibitors)

RN 773882-10-3 CAPLUS

CN Benzamide, 4-[4-(diphenylmethyl)-1-piperaziny]-3-
[[(phenylamino)carbonyl]amino]- (9CI) (CA INDEX NAME)



=> file caold

COST IN U.S. DOLLARS

SINCE FILE

TOTAL

ENTRY

SESSION

FULL ESTIMATED COST

5.74

196.80

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)

SINCE FILE

TOTAL

ENTRY

SESSION

CA SUBSCRIBER PRICE

-0.78

-0.78

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FILE COVERS 1907-1966

FILE LAST UPDATED: 01 May 1997 (19970501/UP)

This file contains CAS Registry Numbers for easy and accurate substance identification. Title keywords, authors, patent assignees, and patent information, e.g., patent numbers, are now searchable from 1907-1966. TIFF images of CA abstracts printed between 1907-1966 are available in the PAGE display formats.

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This file supports REGISTRY for direct browsing and searching of all substance data from the REGISTRY file. Enter HELP FIRST for more information.

=> s 15

L7 0 L5

=> file chemcats

COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	0.45	197.25
DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)	SINCE FILE ENTRY	TOTAL SESSION
CA SUBSCRIBER PRICE	0.00	-0.78

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FILE LAST UPDATED 21 APRIL 2007 (20070421/UP)

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=> s 15

L8 287 L5

=> file reg

COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	0.93	198.18
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CA SUBSCRIBER PRICE	0.00	-0.78

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STRUCTURE FILE UPDATES: 27 APR 2007 HIGHEST RN 933069-51-3
DICTIONARY FILE UPDATES: 27 APR 2007 HIGHEST RN 933069-51-3

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TSCA INFORMATION NOW CURRENT THROUGH December 2, 2006

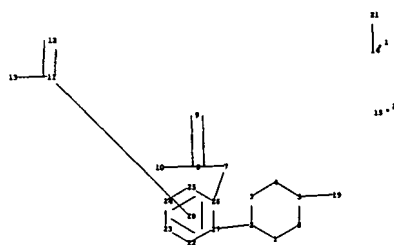
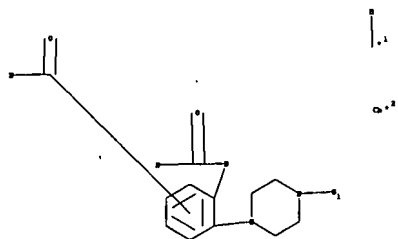
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<http://www.cas.org/support/stngen/stndoc/properties.html>

=>

Uploading C:\Documents and Settings\EBernhardt\My Documents\Stnexp\Queries\10815017-B.str



chain nodes :
7 8 9 10 11 12 14 15 19 21
ring nodes :
1 2 3 4 5 6 22 23 24 25 26 27
ring/chain nodes :

10/815017

13

chain bonds :

2-27 5-19 7-8 7-26 8-9 8-10 11-12 11-13 14-21

ring bonds :

1-2 1-6 2-3 3-4 4-5 5-6 22-23 22-27 23-24 24-25 25-26 26-27

exact/norm bonds :

1-2 1-6 2-3 2-27 3-4 4-5 5-6 5-19 7-8 7-26 8-9 8-10 11-12 11-13

exact bonds :

14-21

normalized bonds :

22-23 22-27 23-24 24-25 25-26 26-27

isolated ring systems :

containing 1 :

G1:[*1],[*2]

Match level :

1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:CLASS 8:CLASS 9:CLASS 10:CLASS
11:CLASS 12:CLASS 13:CLASS 14:CLASS 15:Atom 19:CLASS 21:CLASS 22:CLASS
23:Atom 24:Atom 25:Atom 26:CLASS 27:Atom 28:Atom

L9 STRUCTURE UPLOADED

=> s 19

SAMPLE SEARCH INITIATED 16:17:02 FILE 'REGISTRY'

SAMPLE SCREEN SEARCH COMPLETED - 36 TO ITERATE

100.0% PROCESSED 36 ITERATIONS

1 ANSWERS

SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
BATCH **COMPLETE**

PROJECTED ITERATIONS: 360 TO 1080

PROJECTED ANSWERS: 1 TO 80

L10 1 SEA SSS SAM L9

=> d his

(FILE 'HOME' ENTERED AT 16:07:18 ON 28 APR 2007)

FILE 'REGISTRY' ENTERED AT 16:07:30 ON 28 APR 2007

L1 STRUCTURE UPLOADED

L2 34 S L1

L3 STRUCTURE UPLOADED

L4 27 S L3

L5 422 S L3 SSS FULL

SAVE L5 TEN815017/A

FILE 'CAPLUS' ENTERED AT 16:12:27 ON 28 APR 2007

L6 1 S L5

FILE 'CAOLD' ENTERED AT 16:13:01 ON 28 APR 2007

L7 0 S L5

10/815017

FILE 'CHEMCATS' ENTERED AT 16:13:11 ON 28 APR 2007
L8 287 S L5

FILE 'REGISTRY' ENTERED AT 16:13:45 ON 28 APR 2007
L9 STRUCTURE UPLOADED
L10 1 S L9

=> s l9 sub=15 full
FULL SUBSET SEARCH INITIATED 16:18:43 FILE 'REGISTRY'
FULL SUBSET SCREEN SEARCH COMPLETED - 63 TO ITERATE

100.0% PROCESSED 63 ITERATIONS 63 ANSWERS
SEARCH TIME: 00.00.01

L11 63 SEA SUB=L5 SSS FUL L9

=> file caplus		
COST IN U.S. DOLLARS	SINCE FILE	TOTAL
	ENTRY	SESSION
FULL ESTIMATED COST	44.25	242.43
DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)	SINCE FILE	TOTAL
	ENTRY	SESSION
CA SUBSCRIBER PRICE	0.00	-0.78

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=> s l11
L12 1 L11

=> d l12 bib abs hitstr

L12 ANSWER 1 OF 1 CAPLUS COPYRIGHT 2007 ACS on STN
AN 2004:857583 CAPLUS
DN 141:332220

TI A preparation of (piperazinyphenyl)urea derivatives as phospholipase C inhibitors, useful for the treatment of inflammatory disorders
 IN Lagu, Bharat; Wachter, Michael; Rupert, Kenneth; Wachter, Michael
 PA Janssen Pharmaceutica N.V., Belg.
 SO PCT Int. Appl., 141 pp.
 CODEN: PIXXD2
 DT Patent
 LA English
 FAN.CNT 1

	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
PI	WO 2004087685	A2	20041014	WO 2004-US9846	20040331
	WO 2004087685	A3	20041216		
	W:	AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BW, BY, BZ, CA, CH, CN, CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, EG, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NA, NI, NO, NZ, OM, PG, PH, PL, PT, RO, RU, SC, SD, SE, SG, SK, SL, SY, TJ, TM, TN, TR, TT, TZ, UA, UG, US, UZ, VC, VN, YU, ZA, ZM, ZW			
	RW:	BW, GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZM, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM, AT, BE, BG, CH, CY, CZ, DE, DK, EE, ES, FI, FR, GB, GR, HU, IE, IT, LU, MC, NL, PL, PT, RO, SE, SI, SK, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG			
	US 2004235855	A1	20041125	US 2004-815017	20040331
PRAI	US 2003-458938P	P	20030331		
OS	MARPAT 141:332220				
GI					

* STRUCTURE DIAGRAM TOO LARGE FOR DISPLAY - AVAILABLE VIA OFFLINE PRINT *

AB The invention relates to a preparation of (piperazinyphenyl)urea derivs. of formula I [wherein: X is NH₂, NH-alkyl, NHOH, NH-CN, or heterocyclic ring, etc.; Y is one or more (un)substituted alkyl; Z is (CH₂)₂₋₅; R₁ is (un)substituted alkyl, cycloalkyl, or aryl, etc.; R₂ is (un)substituted alkyl, C(O)alkyl, C(O)alkenyl, aryl, or cycloalkyl, etc.; R₃ is O or S], useful as PLC- β 2 inhibitors. For instance, (piperazinyphenyl)urea derivative II (IC₅₀ = 1.2 μ M) was prepared via addition of resin-bound (piperazinyphenyl)amine derivative III to Ph-N=C=O and subsequent resin cleavage (example 1).

IT 773882-10-3P 773882-11-4P 773882-12-5P
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 773882-16-9P 773882-17-0P 773882-18-1P
 773882-19-2P 773882-20-5P 773882-21-6P
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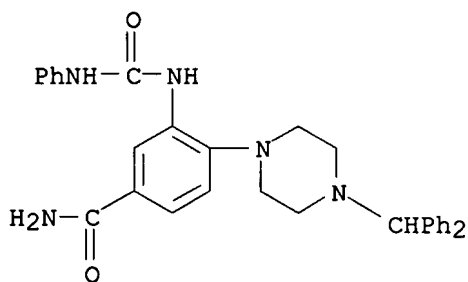
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 773882-86-3P 773882-87-4P 773882-88-5P

RL: PAC (Pharmacological activity); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)

(preparation of (piperazinyphenyl)urea derivs. useful as PLC- β 2 inhibitors)

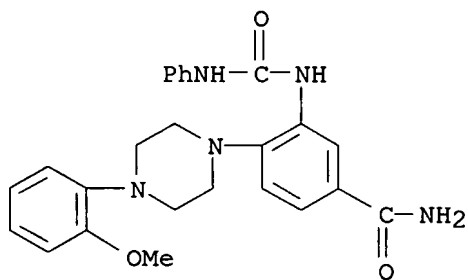
RN 773882-10-3 CAPLUS

CN Benzamide, 4-[4-(diphenylmethyl)-1-piperazinyl]-3-
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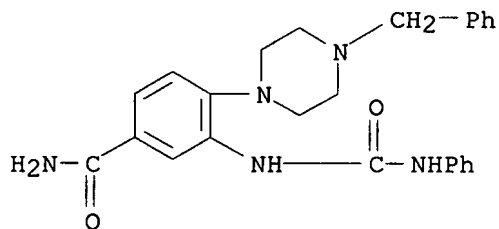
RN 773882-11-4 CAPLUS

CN Benzamide, 4-[4-(2-methoxyphenyl)-1-piperazinyl]-3-
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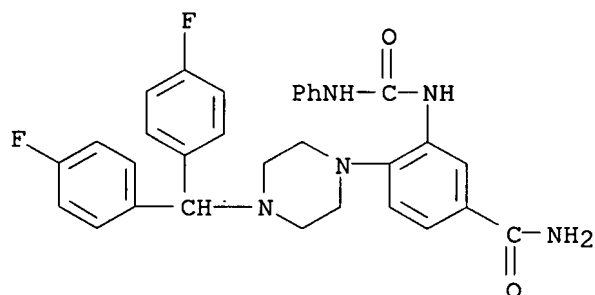
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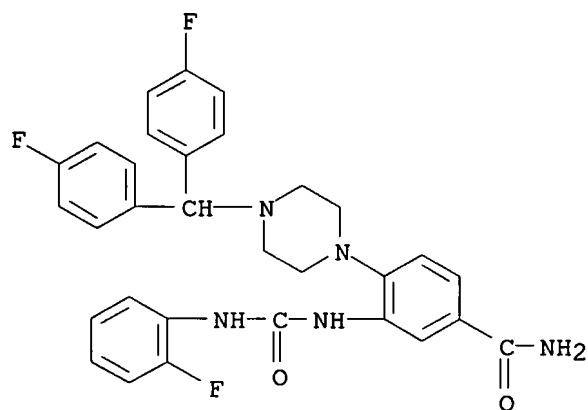
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CN Benzamide, 4-[4-[bis(4-fluorophenyl)methyl]-1-piperazinyl]-3-[[4-[(phenylamino)carbonyl]amino]- (9CI) (CA INDEX NAME)



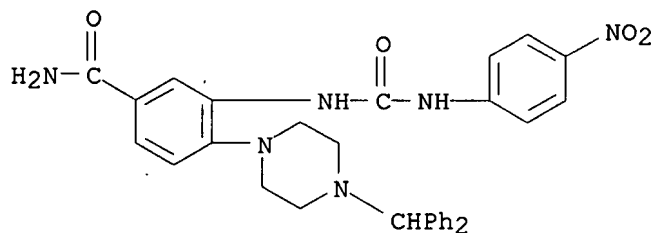
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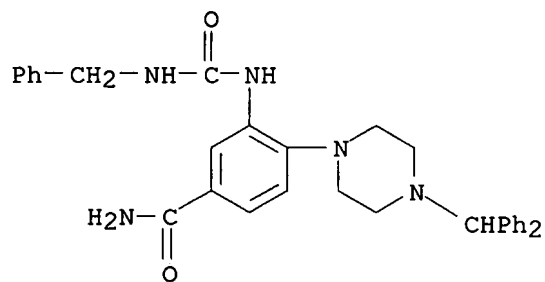


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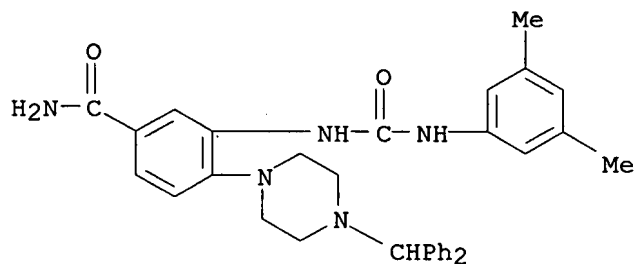


RN 773882-16-9 CAPLUS

CN Benzamide, 4-[4-(diphenylmethyl)-1-piperazinyl]-3-
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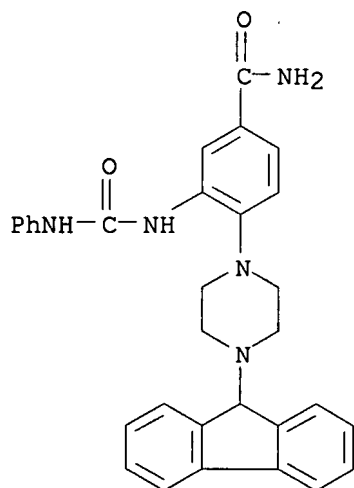
RN 773882-17-0 CAPLUS

CN Benzamide, 3-[[[(3,5-dimethylphenyl)amino]carbonyl]amino]-4-[4-(diphenylmethyl)-1-piperazinyl]- (9CI) (CA INDEX NAME)



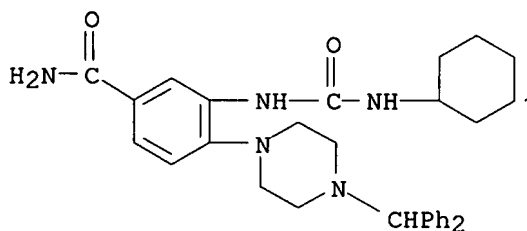
RN 773882-18-1 CAPLUS

CN Benzamide, 4-[4-(9H-fluoren-9-yl)-1-piperazinyl]-3-
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RN 773882-19-2 CAPLUS

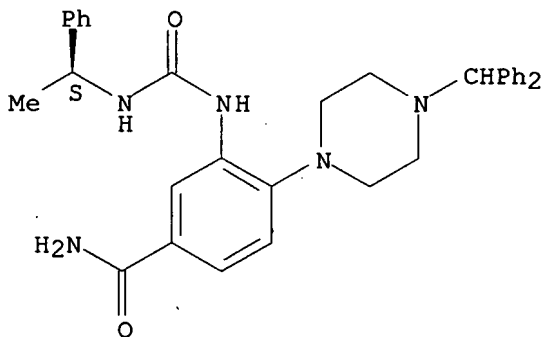
CN Benzamide, 3-[[[(cyclohexylamino)carbonyl]amino]-4-[4-(diphenylmethyl)-1-piperazinyl]]- (9CI) (CA INDEX NAME)



RN 773882-20-5 CAPLUS

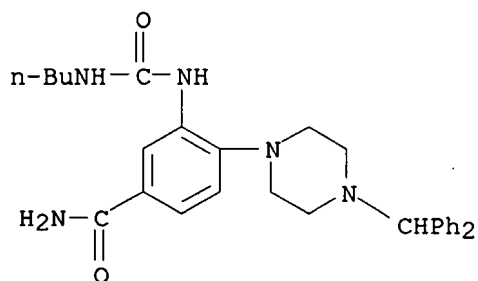
CN Benzamide, 4-[4-(diphenylmethyl)-1-piperazinyl]-3-[[[(1S)-1-phenylethyl]amino]carbonyl]amino]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.



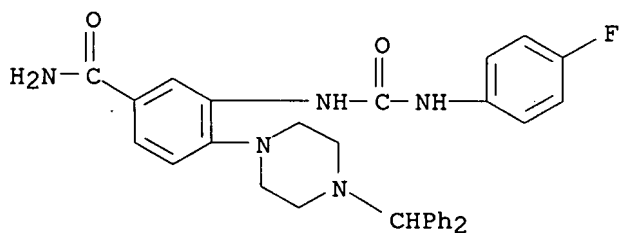
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CN Benzamide, 3-[[[(butylamino)carbonyl]amino]-4-[4-(diphenylmethyl)-1-piperazinyl]]- (9CI) (CA INDEX NAME)



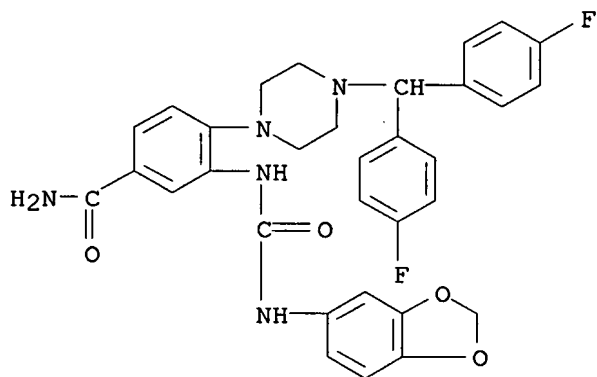
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CN Benzamide, 4-[4-(diphenylmethyl)-1-piperazinyl]-3-[[[(4-fluorophenyl)amino]carbonyl]amino]- (9CI) (CA INDEX NAME)



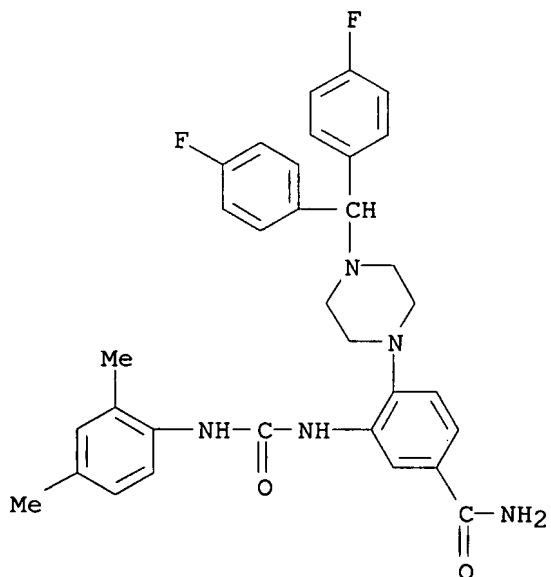
RN 773882-23-8 CAPLUS

CN Benzamide, 3-[[[(1,3-benzodioxol-5-ylamino)carbonyl]amino]-4-[4-[bis(4-fluorophenyl)methyl]-1-piperazinyl]]- (9CI) (CA INDEX NAME)



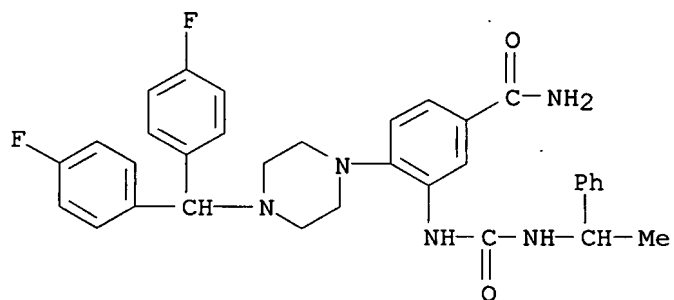
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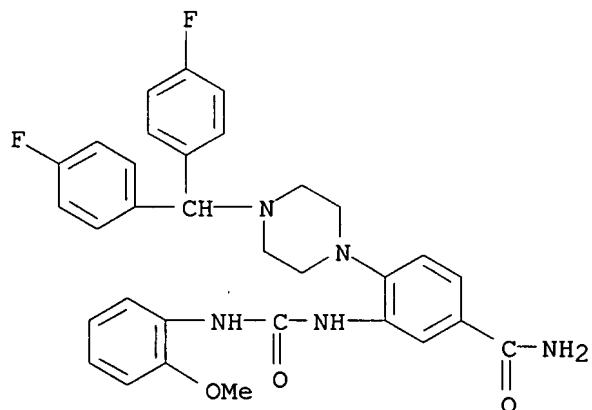
RN 773882-25-0 CAPLUS

CN Benzamide, 4-[4-[bis(4-fluorophenyl)methyl]-1-piperazinyl]-3-[[[(1-phenylethyl)amino]carbonyl]amino]- (9CI) (CA INDEX NAME)



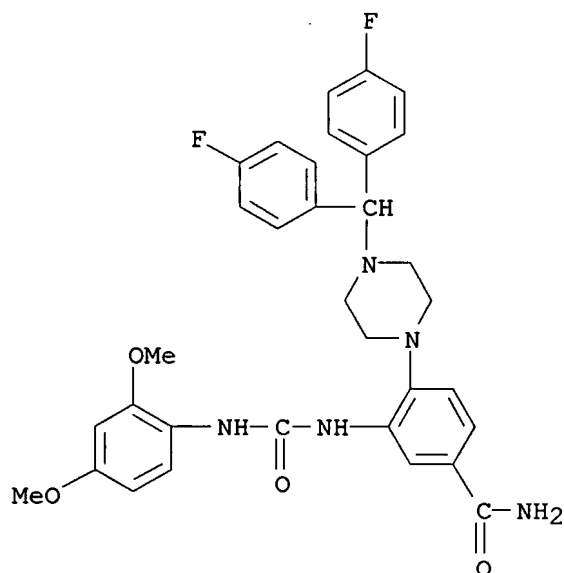
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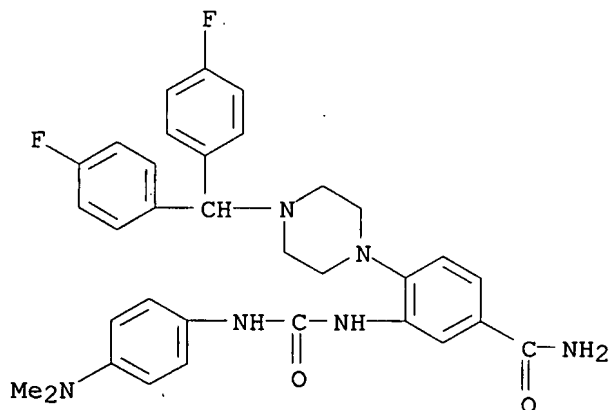
RN 773882-27-2 CAPLUS

CN Benzamide, 4-[4-[bis(4-fluorophenyl)methyl]-1-piperazinyl]-3-[[[(2,4-dimethoxyphenyl)amino]carbonyl]amino]- (9CI) (CA INDEX NAME)



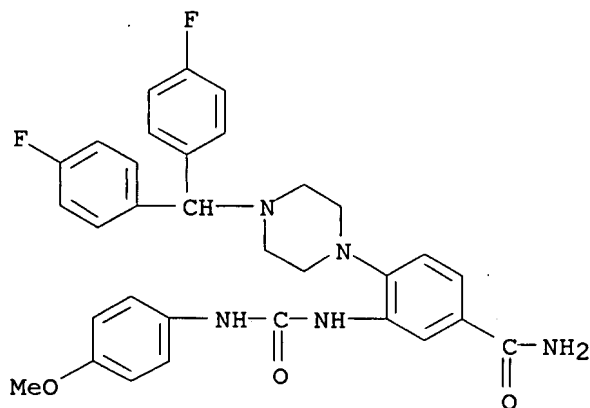
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CN Benzamide, 4-[4-[bis(4-fluorophenyl)methyl]-1-piperazinyl]-3-[[[4-(dimethylamino)phenyl]amino]carbonyl]amino]- (9CI) (CA INDEX NAME)



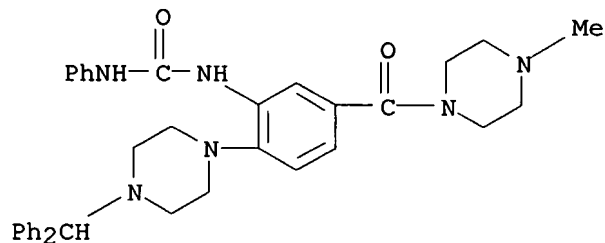
RN 773882-29-4 CAPLUS

CN Benzamide, 4-[4-[bis(4-fluorophenyl)methyl]-1-piperazinyl]-3-[[[(4-methoxyphenyl)amino]carbonyl]amino]- (9CI) (CA INDEX NAME)



RN 773882-32-9 CAPLUS

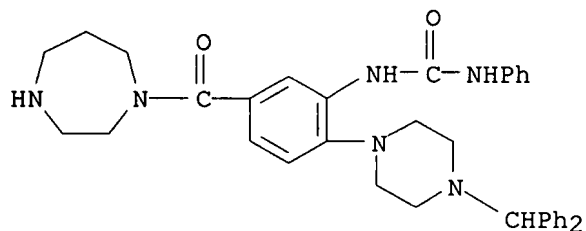
CN Piperazine, 1-[4-[4-(diphenylmethyl)-1-piperazinyl]-3-[[[(phenylamino)carbonyl]amino]benzoyl]-4-methyl- (9CI) (CA INDEX NAME)



RN 773882-33-0 CAPLUS

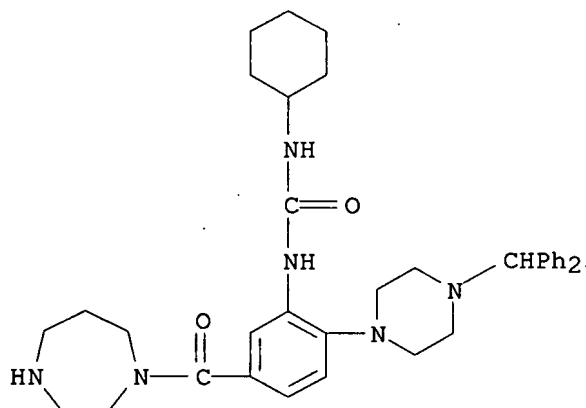
CN 1H-1,4-Diazepine, 1-[4-[4-(diphenylmethyl)-1-piperazinyl]-3-[[[(phenylamino)carbonyl]amino]benzoyl]hexahydro- (9CI) (CA INDEX NAME)

10/815017



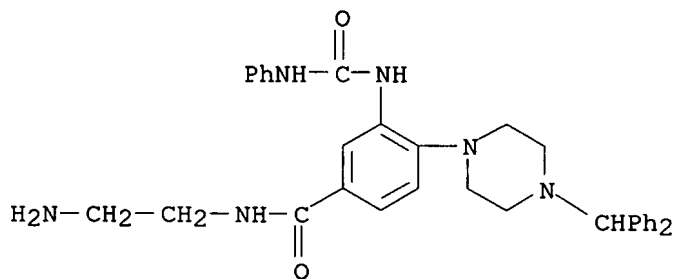
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CN 1H-1,4-Diazepine, 1-[3-[[[(cyclohexylamino)carbonyl]amino]-4-[4-(diphenylmethyl)-1-piperazinyl]benzoyl]hexahydro- (9CI) (CA INDEX NAME)



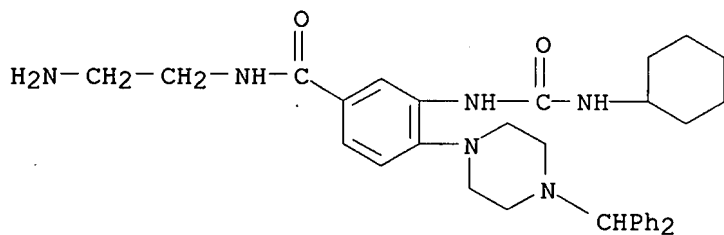
RN 773882-35-2 CAPLUS

CN Benzamide, N-(2-aminoethyl)-4-[4-(diphenylmethyl)-1-piperazinyl]-3-[[[(phenylamino)carbonyl]amino]- (9CI) (CA INDEX NAME)



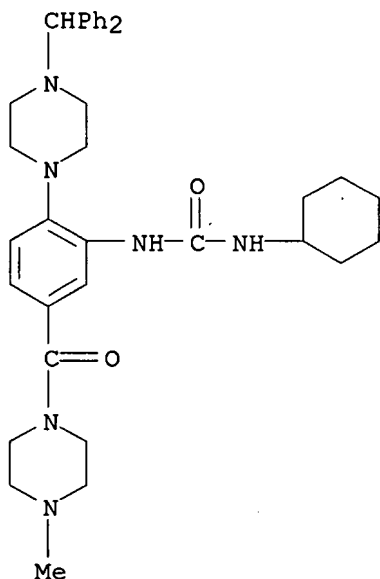
RN 773882-36-3 CAPLUS

CN Benzamide, N-(2-aminoethyl)-3-[[[(cyclohexylamino)carbonyl]amino]-4-[4-(diphenylmethyl)-1-piperazinyl]- (9CI) (CA INDEX NAME)



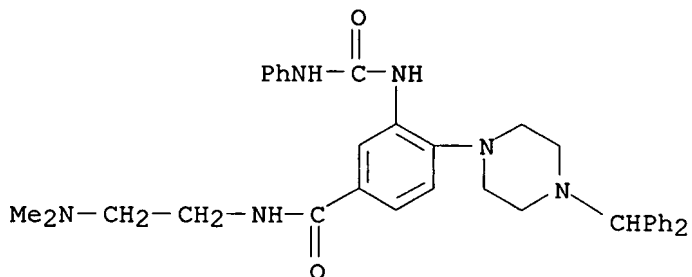
RN 773882-37-4 CAPLUS

CN Piperazine, 1-[3-[[[(cyclohexylamino)carbonyl]amino]-4-[4-(diphenylmethyl)-1-piperazinyl]benzoyl]-4-methyl- (9CI) (CA INDEX NAME)



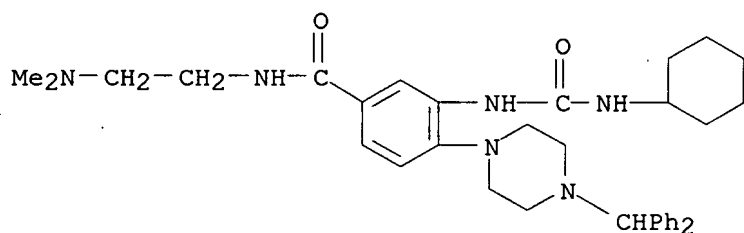
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CN Benzamide, N-[2-(dimethylamino)ethyl]-4-[4-(diphenylmethyl)-1-piperazinyl]-3-[[[(phenylamino)carbonyl]amino]- (9CI) (CA INDEX NAME)



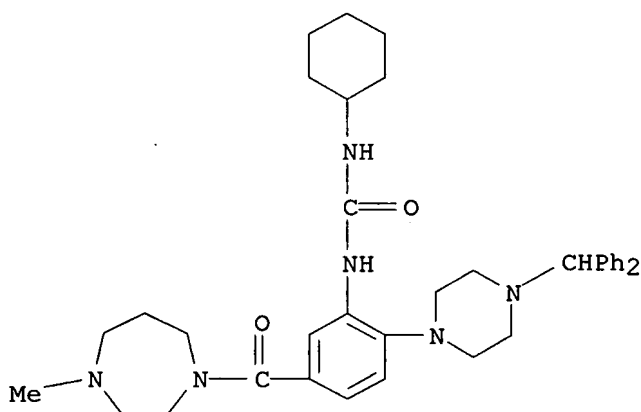
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CN Benzamide, 3-[[[(cyclohexylamino)carbonyl]amino]-N-[2-(dimethylamino)ethyl]-4-[4-(diphenylmethyl)-1-piperazinyl]- (9CI) (CA INDEX NAME)



RN 773882-40-9 CAPLUS

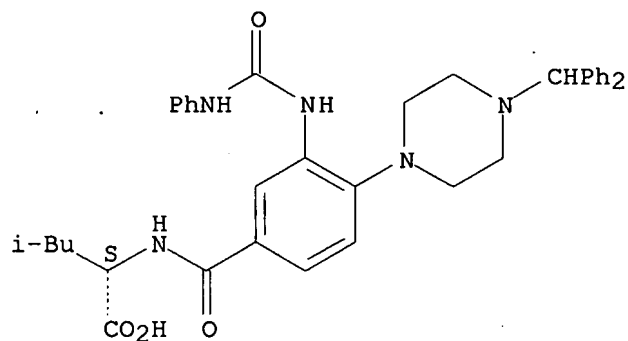
CN 1H-1,4-Diazepine, 1-[3-[[(cyclohexylamino) carbonyl] amino]-4-[4-(diphenylmethyl)-1-piperazinyl]benzoyl]hexahydro-4-methyl- (9CI) (CA INDEX NAME)



RN 773882-41-0 CAPLUS

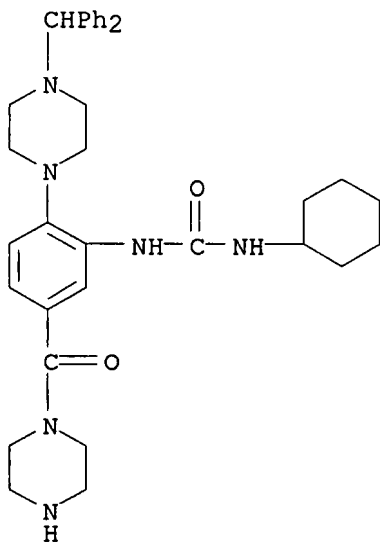
CN L-Leucine, N-[4-[4-(diphenylmethyl)-1-piperazinyl]-3-[[(phenylamino) carbonyl] amino]benzoyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.



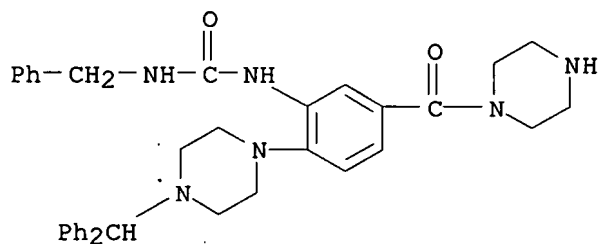
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CN Piperazine, 1-[3-[[(cyclohexylamino) carbonyl] amino]-4-[4-(diphenylmethyl)-1-piperazinyl]benzoyl]- (9CI) (CA INDEX NAME)



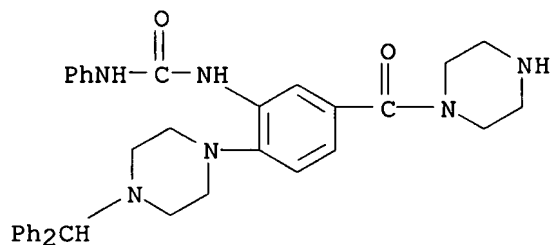
RN 773882-43-2 CAPLUS

CN Piperazine, 1-[4-[4-(diphenylmethyl)-1-piperazinyl]-3-
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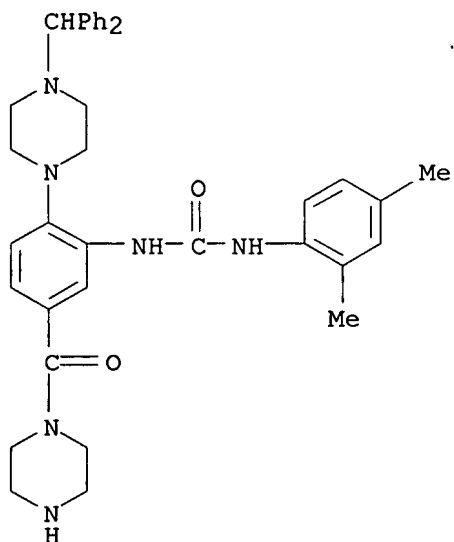
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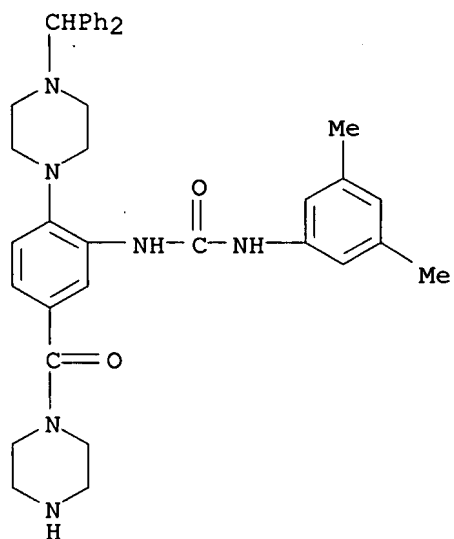
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CN Piperazine, 1-[3-[[[(2,4-dimethylphenyl)amino]carbonyl]amino]-4-[4-
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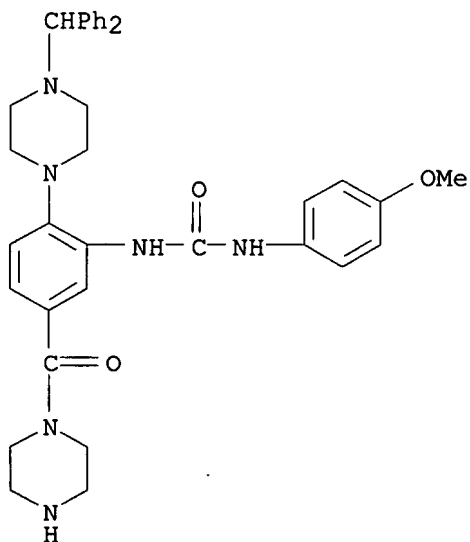
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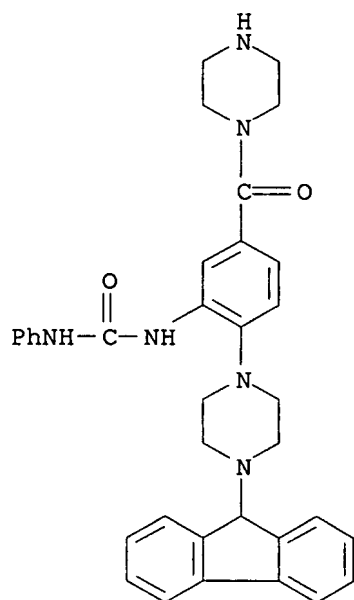
CN Piperazine, 1-[4-[4-(diphenylmethyl)-1-piperazinyl]-3-[[[(4-methoxyphenyl)amino]carbonyl]amino]benzoyl]- (9CI) (CA INDEX NAME)

10/815017



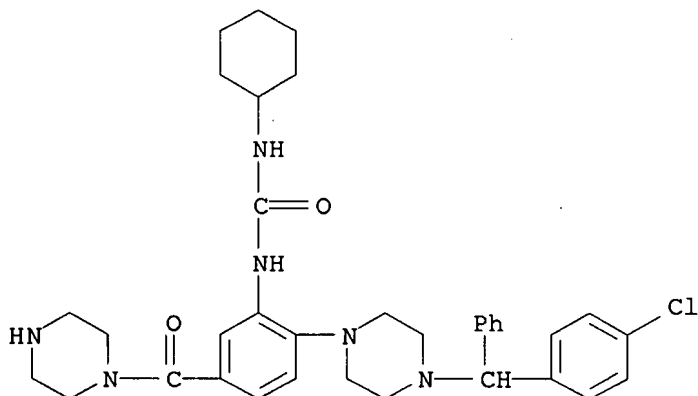
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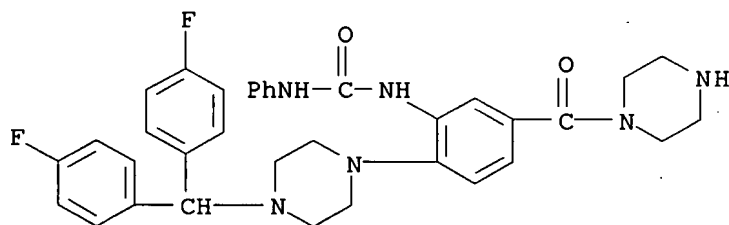
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CN Piperazine, 1-[4-[4-[(4-chlorophenyl)phenylmethyl]-1-piperazinyl]-3-
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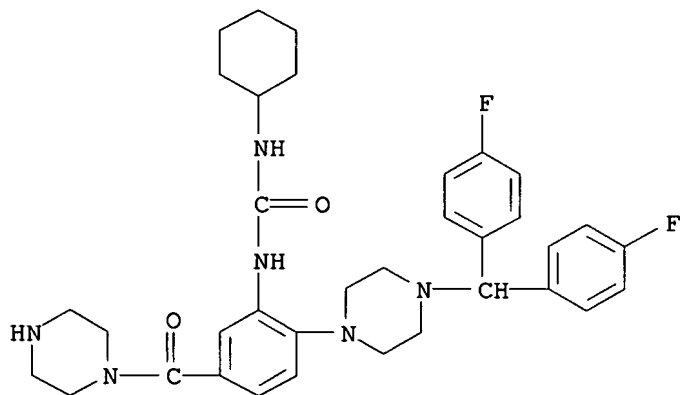
RN 773882-50-1 CAPLUS

CN Piperazine, 1-[4-[4-[(cyclohexylamino)carbonyl]amino]benzoyl]-3-[[4-(4-chlorophenyl)-1-phenylethyl]-1-piperazinyl]-3-[[4-(4-chlorophenyl)-1-phenylethyl]-1-piperazinyl] (9CI) (CA INDEX NAME)



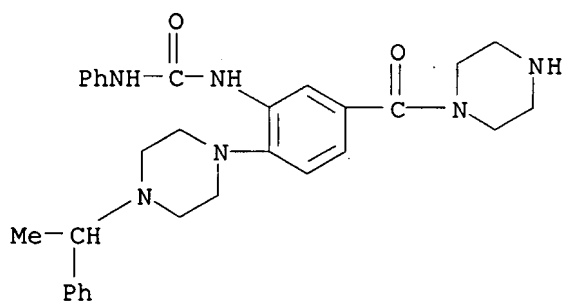
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CN Piperazine, 1-[4-[4-[(bis(4-fluorophenyl)methyl)amino]benzoyl]-1-piperazinyl]-3-[[4-(4-fluorophenyl)-1-phenylethyl]-1-piperazinyl]-3-[[4-(4-fluorophenyl)-1-phenylethyl]-1-piperazinyl] (9CI) (CA INDEX NAME)

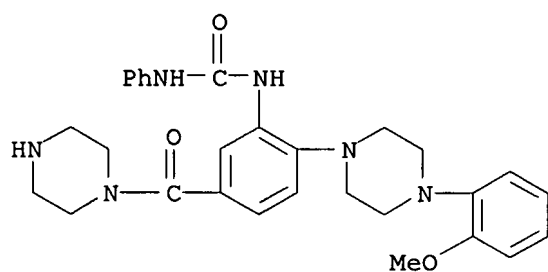


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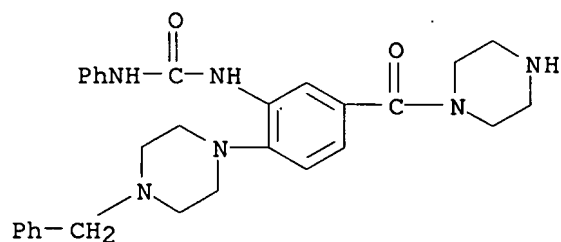
CN Piperazine, 1-[3-[[4-(1-phenylethyl)-1-piperazinyl]benzoyl]-4-[[4-(1-phenylethyl)-1-piperazinyl]benzoyl]-4-[[4-(1-phenylethyl)-1-piperazinyl]benzoyl]-4-[[4-(1-phenylethyl)-1-piperazinyl]benzoyl] (9CI) (CA INDEX NAME)



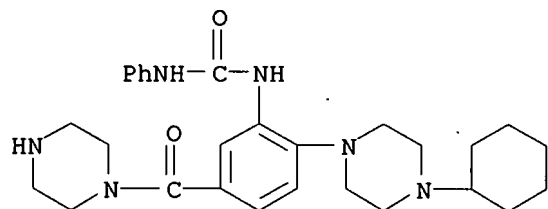
RN 773882-56-7 CAPLUS

CN Piperazine, 1-[4-[4-(2-methoxyphenyl)-1-piperazinyl]-3-
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RN 773882-58-9 CAPLUS

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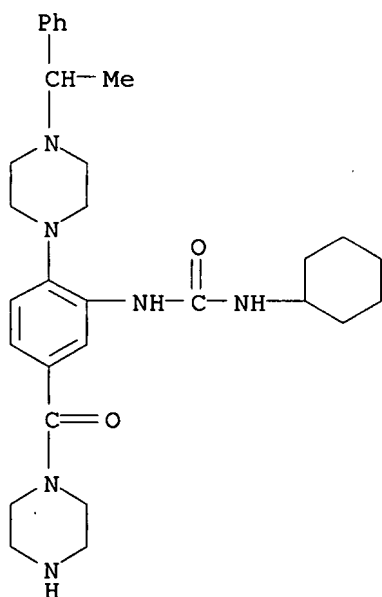
RN 773882-59-0 CAPLUS

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[[(phenylamino)carbonyl]amino]benzoyl]- (9CI) (CA INDEX NAME)

10/815017

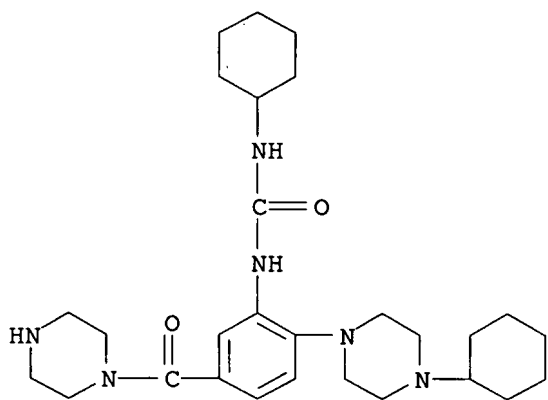
RN 773882-61-4 CAPLUS

CN Piperazine, 1-[3-[[[(cyclohexylamino)carbonyl]amino]-4-[4-(1-phenylethyl)-1-piperazinyl]benzoyl]- (9CI) (CA INDEX NAME)



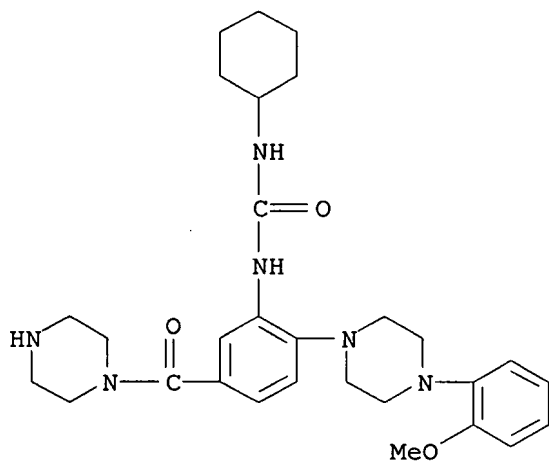
RN 773882-63-6 CAPLUS

CN Piperazine, 1-[3-[[[(cyclohexylamino)carbonyl]amino]-4-(4-cyclohexyl-1-piperazinyl)benzoyl]- (9CI) (CA INDEX NAME)



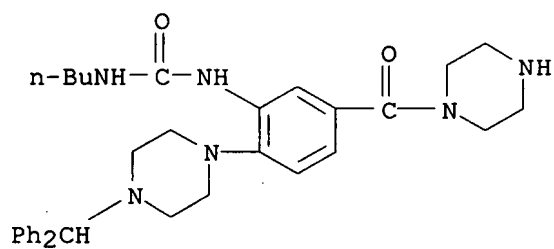
RN 773882-65-8 CAPLUS

CN Piperazine, 1-[3-[[[(cyclohexylamino)carbonyl]amino]-4-[4-(2-methoxyphenyl)-1-piperazinyl]benzoyl]- (9CI) (CA INDEX NAME)



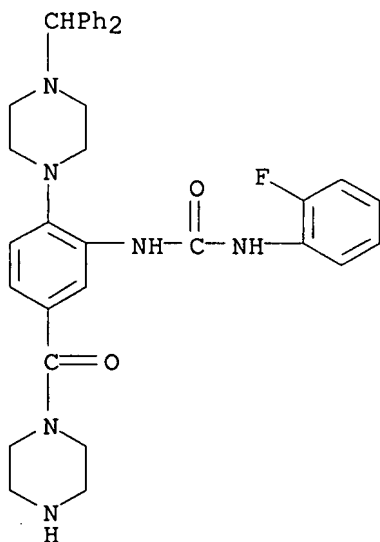
RN 773882-66-9 CAPLUS

CN Piperazine, 1-[3-[[4-(butylamino)carbonyl]amino]-4-[4-(diphenylmethyl)-1-piperazinyl]benzoyl]- (9CI) (CA INDEX NAME)



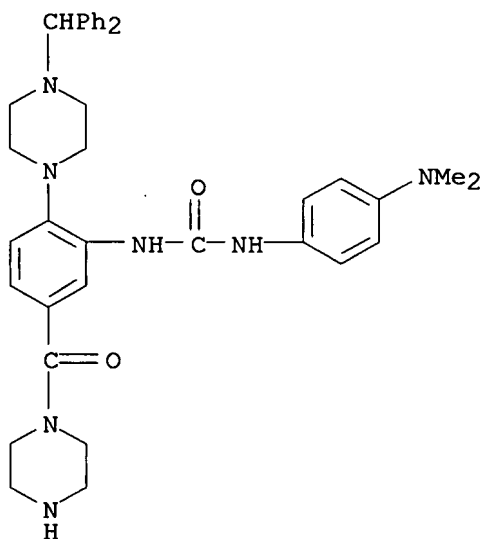
RN 773882-67-0 CAPLUS

CN Piperazine, 1-[4-[4-(diphenylmethyl)-1-piperazinyl]-3-[[[2-(2-fluorophenyl)amino]carbonyl]amino]benzoyl]- (9CI) (CA INDEX NAME)



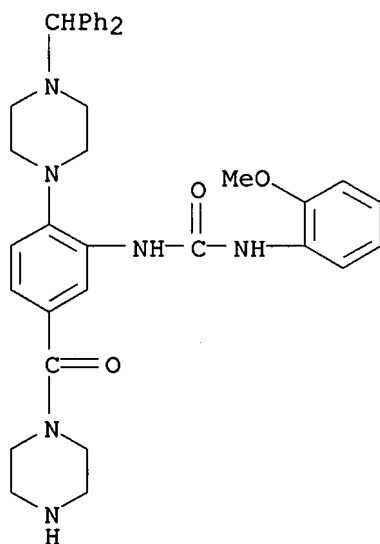
RN 773882-68-1 CAPLUS

CN Piperazine, 1-[3-[[[4-(dimethylamino)phenyl]amino]carbonyl]amino]-4-[4-(diphenylmethyl)-1-piperazinyl]benzoyl]- (9CI) (CA INDEX NAME)

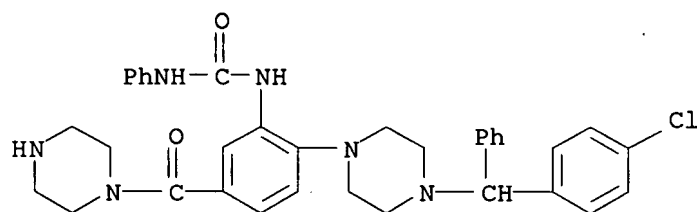


RN 773882-69-2 CAPLUS

CN Piperazine, 1-[4-[4-(diphenylmethyl)-1-piperazinyl]-3-[[[(2-methoxyphenyl)amino]carbonyl]amino]benzoyl]- (9CI) (CA INDEX NAME)



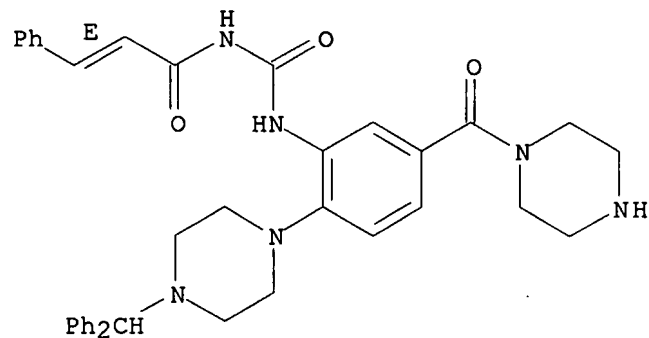
RN 773882-70-5 CAPLUS

CN Piperazine, 1-[4-[4-[(4-chlorophenyl)phenylmethyl]-1-piperazinyl]-3-
[[(phenylamino) carbonyl] amino] benzoyl]- (9CI) (CA INDEX NAME)

RN 773882-71-6 CAPLUS

CN 2-Propenamide, N-[[[2-[4-(diphenylmethyl)-1-piperazinyl]-5-(1-
piperazinylcarbonyl)phenyl]amino]carbonyl]-3-phenyl-, (2E)- (9CI) (CA
INDEX NAME)

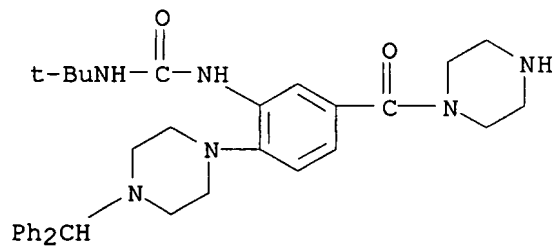
Double bond geometry as shown.



RN 773882-72-7 CAPLUS

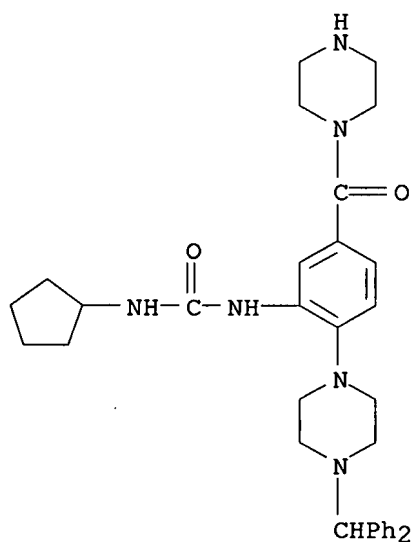
10/815017

CN Piperazine, 1-[3-[[[(1,1-dimethylethyl)amino]carbonyl]amino]-4-[4-(diphenylmethyl)-1-piperazinyl]benzoyl]- (9CI) (CA INDEX NAME)



RN 773882-73-8 CAPLUS

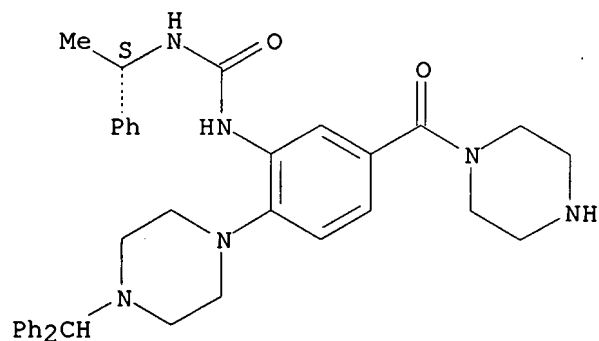
CN Piperazine, 1-[3-[[[(cyclopentylamino)carbonyl]amino]-4-[4-(diphenylmethyl)-1-piperazinyl]benzoyl]- (9CI) (CA INDEX NAME)



RN 773882-74-9 CAPLUS

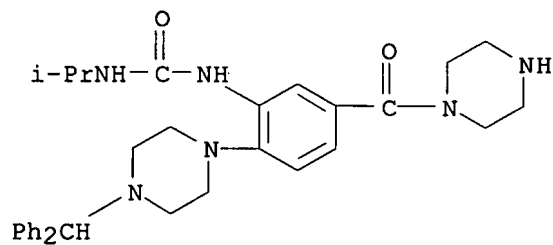
CN Piperazine, 1-[4-[4-(diphenylmethyl)-1-piperazinyl]-3-[[[(1S)-1-phenylethyl]amino]carbonyl]amino]benzoyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.



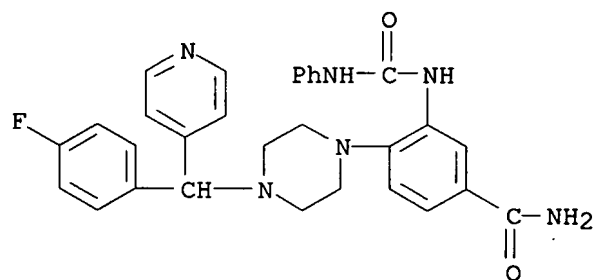
RN 773882-76-1 CAPLUS

CN Piperazine, 1-[4-[4-(diphenylmethyl)-1-piperazinyl]-3-[[[(1-methylethyl)amino]carbonyl]amino]benzoyl]- (9CI) (CA INDEX NAME)



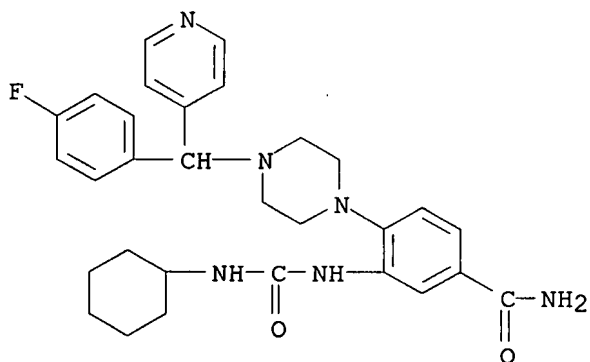
RN 773882-78-3 CAPLUS

CN Benzamide, 4-[4-[(4-fluorophenyl)-4-pyridinylmethyl]-1-piperazinyl]-3-[[[(phenylamino)carbonyl]amino]- (9CI) (CA INDEX NAME)



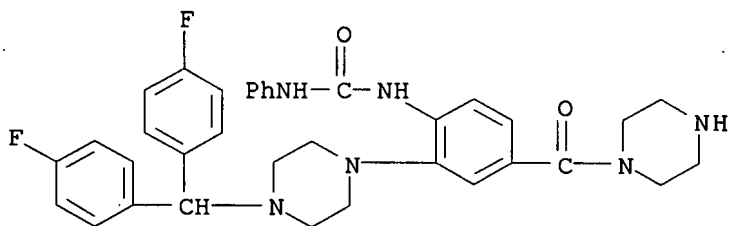
RN 773882-79-4 CAPLUS

CN Benzamide, 3-[[[(cyclohexylamino)carbonyl]amino]-4-[4-[(4-fluorophenyl)-4-pyridinylmethyl]-1-piperazinyl]- (9CI) (CA INDEX NAME)



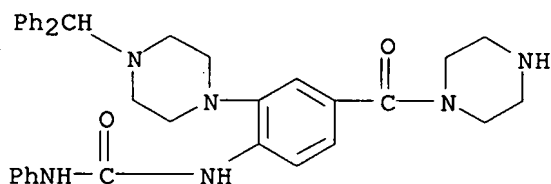
RN 773882-85-2 CAPLUS

CN Piperazine, 1-[3-[4-[bis(4-fluorophenyl)methyl]-1-piperazinyl]-4-[[phenylamino)carbonyl]amino]benzoyl]- (9CI) (CA INDEX NAME)



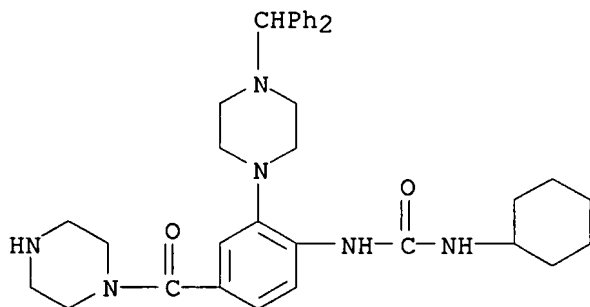
RN 773882-86-3 CAPLUS

CN Piperazine, 1-[3-[4-(diphenylmethyl)-1-piperazinyl]-4-[[phenylamino)carbonyl]amino]benzoyl]- (9CI) (CA INDEX NAME)

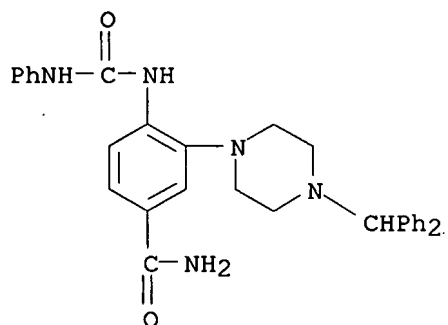


RN 773882-87-4 CAPLUS

CN Piperazine, 1-[4-[[cyclohexylamino)carbonyl]amino]-3-[4-(diphenylmethyl)-1-piperazinyl]benzoyl]- (9CI) (CA INDEX NAME)



RN 773882-88-5 CAPLUS
 CN Benzamide, 3-[4-(diphenylmethyl)-1-piperazinyl]-4-
 [[(phenylamino)carbonyl]amino]- (9CI) (CA INDEX NAME)



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FULL ESTIMATED COST

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SINCE FILE	TOTAL
ENTRY	SESSION
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